



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2024 – 09:24 AM EST

PDB ID : 2Z5H  
Title : Crystal structure of the head-to-tail junction of tropomyosin complexed with a fragment of TnT  
Authors : Murakami, K.; Nozawa, K.; Tomii, K.; Kudou, N.; Igarashi, N.; Shirakihara, Y.; Wakatsuki, S.; Stewart, M.; Yasunaga, T.; Wakabayashi, T.  
Deposited on : 2007-07-12  
Resolution : 2.89 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

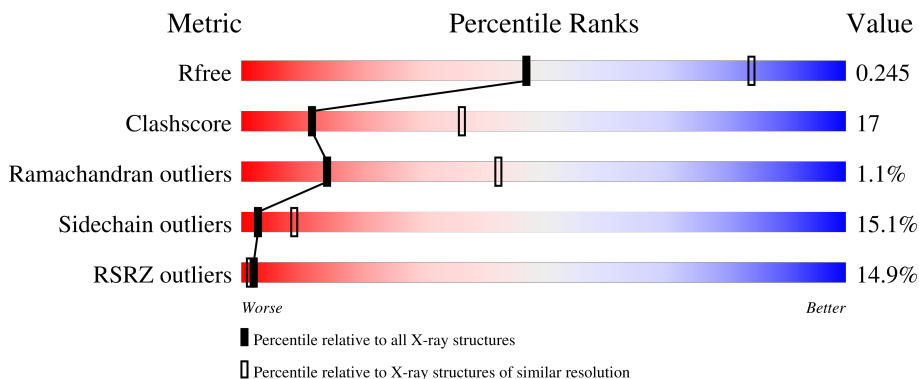
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.89 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	52	 2% 65% 25% 6%
1	B	52	 2% 83% 12%
1	C	52	 2% 75% 21%
1	D	52	 6% 83% 12% 6%
1	E	52	 8% 83% 8% 6%

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Mol	Chain	Length	Quality of chain
1	F	52	<p>4% 81% 17% .</p>
1	G	52	<p>4% 69% 27% . .</p>
1	H	52	<p>2% 81% 12% . 6%</p>
2	I	40	<p>30% 90% 48% 20% .</p>
3	T	55	<p>5% 44% 33% 18% 5% 38%</p>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3940 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called General control protein GCN4 and Tropomyosin alpha-1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	49	Total 403	C 254	N 66	O 82	S 1	0	0	0
1	B	51	Total 416	C 261	N 68	O 86	S 1	0	0	0
1	C	52	Total 424	C 267	N 69	O 87	S 1	0	0	0
1	D	49	Total 403	C 254	N 66	O 82	S 1	0	0	0
1	E	49	Total 403	C 254	N 66	O 82	S 1	0	0	0
1	F	51	Total 416	C 261	N 68	O 86	S 1	0	0	0
1	G	52	Total 424	C 267	N 69	O 87	S 1	0	0	0
1	H	49	Total 403	C 254	N 66	O 82	S 1	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	233	GLY	-	expression tag	UNP P03069
B	233	GLY	-	expression tag	UNP P03069
C	233	GLY	-	expression tag	UNP P03069
D	233	GLY	-	expression tag	UNP P03069
E	233	GLY	-	expression tag	UNP P03069
F	233	GLY	-	expression tag	UNP P03069
G	233	GLY	-	expression tag	UNP P03069
H	233	GLY	-	expression tag	UNP P03069

- Molecule 2 is a protein called Tropomyosin alpha-1 chain and General control protein GCN4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	39	Total	C	N	O	S	0	0	0
			306	189	56	58	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	-3	GLY	-	expression tag	UNP P58772
I	-2	ALA	-	expression tag	UNP P58772
I	-1	ALA	-	expression tag	UNP P58772
I	0	SER	-	expression tag	UNP P58772

- Molecule 3 is a protein called Troponin T, fast skeletal muscle isoforms.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	T	34	Total	C	N	O	8	0	0
			287	179	53	55			

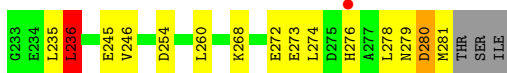
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	7	Total	O	0	0
			7	7		
4	C	10	Total	O	0	0
			10	10		
4	D	3	Total	O	0	0
			3	3		
4	E	4	Total	O	0	0
			4	4		
4	F	12	Total	O	0	0
			12	12		
4	G	2	Total	O	0	0
			2	2		
4	H	2	Total	O	0	0
			2	2		
4	I	7	Total	O	0	0
			7	7		

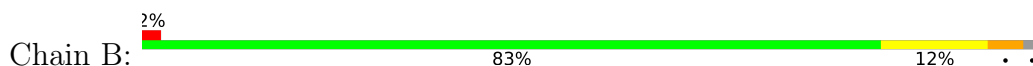
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

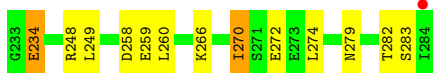
- Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



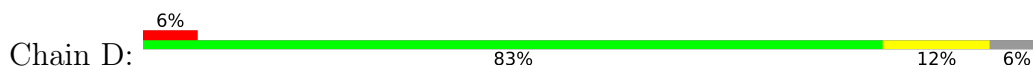
- Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



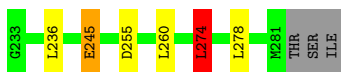
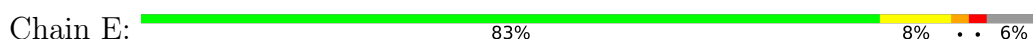
- Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



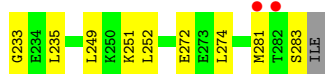
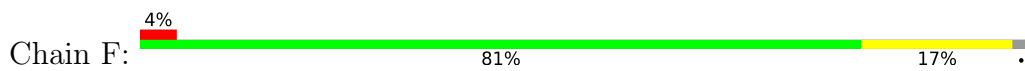
- Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



- Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



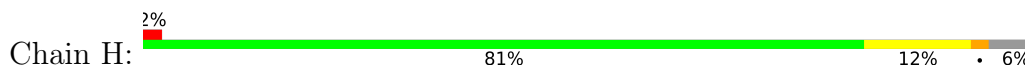
- Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



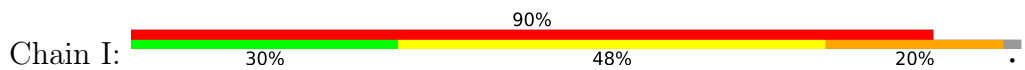
• Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



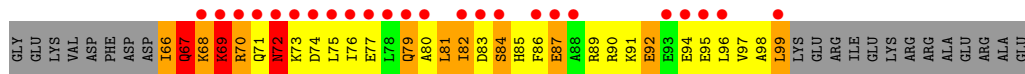
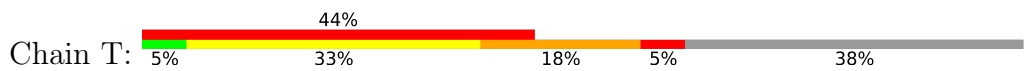
• Molecule 1: General control protein GCN4 and Tropomyosin alpha-1 chain



• Molecule 2: Tropomyosin alpha-1 chain and General control protein GCN4



• Molecule 3: Troponin T, fast skeletal muscle isoforms



## 4 Data and refinement statistics i

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.81Å 158.31Å 163.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	113.96 – 2.89 40.55 – 2.89	Depositor EDS
% Data completeness (in resolution range)	95.9 (113.96-2.89) 95.9 (40.55-2.89)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.16 (at 2.90Å)	Xtrriage
Refinement program	REFMAC 5.2.0019, CNS 1.1	Depositor
R, $R_{free}$	0.237 , 0.247 0.227 , 0.245	Depositor DCC
$R_{free}$ test set	1251 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.9	Xtrriage
Anisotropy	0.730	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 46.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.064 for -h,-l,-k	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3940	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.55	0/407	0.98	2/545 (0.4%)
1	B	0.57	0/420	1.07	0/563
1	C	0.45	0/428	0.83	2/574 (0.3%)
1	D	0.47	0/407	0.81	0/545
1	E	0.51	0/407	0.94	3/545 (0.6%)
1	F	0.51	0/420	0.89	0/563
1	G	0.54	0/428	0.87	0/574
1	H	0.52	0/407	0.81	0/545
2	I	0.40	0/305	0.95	1/401 (0.2%)
3	T	0.76	2/288 (0.7%)	1.37	4/382 (1.0%)
All	All	0.53	2/3917 (0.1%)	0.95	12/5237 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	G	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	T	69	LYS	CB-CG	-8.01	1.30	1.52
3	T	66	ILE	CB-CG2	-6.66	1.32	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	69	LYS	CA-CB-CG	7.17	129.17	113.40
1	E	255	ASP	CB-CG-OD1	6.12	123.81	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	82	ILE	CB-CA-C	-5.64	100.31	111.60
2	I	35	LEU	CA-CB-CG	5.60	128.17	115.30
1	C	248	ARG	NE-CZ-NH1	5.50	123.05	120.30
3	T	69	LYS	CB-CG-CD	-5.47	97.37	111.60
1	E	255	ASP	CB-CG-OD2	-5.34	113.50	118.30
1	A	236	LEU	CB-CG-CD1	-5.30	101.98	111.00
1	C	258	ASP	CB-CG-OD1	5.30	123.07	118.30
3	T	66	ILE	CG1-CB-CG2	5.16	122.76	111.40
1	A	254	ASP	CB-CG-OD1	5.11	122.90	118.30
1	E	274	LEU	CA-CB-CG	5.02	126.85	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	234	GLU	Peptide
1	G	234	GLU	Peptide

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	403	0	400	24	0
1	B	416	0	412	10	0
1	C	424	0	423	3	0
1	D	403	0	400	2	0
1	E	403	0	400	3	0
1	F	416	0	412	3	0
1	G	424	0	423	10	0
1	H	403	0	400	7	0
2	I	306	0	334	29	2
3	T	287	0	297	75	2
4	A	8	0	0	0	0
4	B	7	0	0	1	0
4	C	10	0	0	0	0
4	D	3	0	0	0	0
4	E	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	12	0	0	0	0
4	G	2	0	0	0	0
4	H	2	0	0	0	0
4	I	7	0	0	1	0
All	All	3940	0	3901	134	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (134) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:HIS:CE1	3:T:83:ASP:HB2	1.19	1.67
1:A:276:HIS:CE1	3:T:83:ASP:CB	1.97	1.47
1:A:276:HIS:ND1	3:T:83:ASP:HB2	1.22	1.41
3:T:86:PHE:CE1	3:T:87:GLU:HB2	1.61	1.33
3:T:84:SER:HA	3:T:86:PHE:CE2	1.69	1.25
1:A:276:HIS:HE1	3:T:83:ASP:OD1	1.21	1.19
1:A:276:HIS:ND1	3:T:83:ASP:CB	2.04	1.07
1:A:276:HIS:HE1	3:T:83:ASP:CG	1.58	1.06
1:A:276:HIS:CE1	3:T:83:ASP:OD1	2.12	1.02
3:T:68:LYS:O	3:T:68:LYS:HD2	1.60	1.02
3:T:69:LYS:O	3:T:72:ASN:N	1.94	1.00
2:I:28:GLU:O	2:I:32:LEU:HB2	1.61	1.00
3:T:79:GLN:O	3:T:83:ASP:OD2	1.80	0.97
3:T:92:GLU:O	3:T:96:LEU:HG	1.64	0.97
3:T:86:PHE:CD1	3:T:87:GLU:HB2	2.01	0.96
2:I:21:ARG:O	2:I:25:LEU:HD23	1.67	0.94
3:T:84:SER:HA	3:T:86:PHE:HE2	1.30	0.94
1:A:276:HIS:CE1	3:T:83:ASP:CG	2.34	0.93
2:I:20:ASP:O	2:I:23:GLU:HG2	1.66	0.93
3:T:89:ARG:O	3:T:92:GLU:CG	2.16	0.93
3:T:89:ARG:O	3:T:92:GLU:HG2	1.69	0.90
3:T:84:SER:CA	3:T:86:PHE:CE2	2.56	0.88
3:T:69:LYS:C	3:T:71:GLN:H	1.77	0.87
3:T:86:PHE:CE1	3:T:87:GLU:CB	2.54	0.85
3:T:84:SER:HA	3:T:86:PHE:CD2	2.10	0.84
3:T:81:LEU:HD23	3:T:84:SER:CB	2.06	0.84
3:T:76:ILE:O	3:T:80:ALA:N	2.10	0.84
2:I:20:ASP:HA	2:I:23:GLU:OE2	1.77	0.83
3:T:74:ASP:HA	3:T:77:GLU:HB2	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:291:HOH:O	2:I:1:ALA:HB1	1.82	0.79
3:T:69:LYS:O	3:T:71:GLN:N	2.15	0.79
3:T:84:SER:CA	3:T:86:PHE:CD2	2.67	0.77
3:T:89:ARG:C	3:T:92:GLU:HG2	2.07	0.75
1:B:278:LEU:CD1	2:I:4:ILE:HD12	2.17	0.74
2:I:12:LYS:O	2:I:16:GLU:HG3	1.88	0.74
1:B:281:MET:HA	1:B:281:MET:CE	2.17	0.74
2:I:19:LEU:HA	2:I:22:ALA:HB3	1.68	0.74
3:T:81:LEU:CD2	3:T:84:SER:CB	2.66	0.73
3:T:84:SER:N	3:T:86:PHE:HD2	1.87	0.72
2:I:12:LYS:O	2:I:16:GLU:CG	2.37	0.72
3:T:84:SER:N	3:T:86:PHE:CD2	2.59	0.71
1:A:279:ASN:C	1:A:281:MET:H	1.95	0.71
1:A:279:ASN:HD21	3:T:83:ASP:HB3	1.57	0.69
1:H:234:GLU:H	1:H:234:GLU:CD	1.95	0.69
3:T:89:ARG:O	3:T:92:GLU:HG3	1.94	0.68
3:T:81:LEU:CD2	3:T:84:SER:HB2	2.25	0.67
1:A:280:ASP:HB3	2:I:11:LEU:HD22	1.76	0.67
3:T:86:PHE:HE1	3:T:87:GLU:HB2	1.51	0.67
3:T:73:LYS:O	3:T:77:GLU:HB2	1.93	0.67
3:T:76:ILE:O	3:T:79:GLN:HG3	1.96	0.66
2:I:1:MET:O	4:I:39:HOH:O	2.12	0.66
3:T:83:ASP:C	3:T:86:PHE:CD2	2.69	0.66
1:B:278:LEU:HD12	2:I:4:ILE:HD12	1.79	0.64
3:T:69:LYS:C	3:T:71:GLN:N	2.49	0.64
2:I:15:LYS:HG3	2:I:16:GLU:N	2.13	0.64
3:T:81:LEU:CG	3:T:84:SER:HB2	2.28	0.63
3:T:89:ARG:HG3	3:T:92:GLU:OE2	2.00	0.61
3:T:81:LEU:HD23	3:T:84:SER:HB3	1.81	0.61
1:G:266:LYS:HD3	1:H:267:TYR:CE1	2.35	0.61
3:T:68:LYS:HD2	3:T:68:LYS:C	2.22	0.60
3:T:81:LEU:HG	3:T:84:SER:HB2	1.81	0.60
3:T:89:ARG:CA	3:T:92:GLU:HG2	2.32	0.60
3:T:89:ARG:HA	3:T:92:GLU:HG2	1.84	0.59
3:T:81:LEU:HD23	3:T:84:SER:HB2	1.80	0.59
1:A:276:HIS:ND1	3:T:83:ASP:HB3	2.14	0.58
1:B:281:MET:HA	1:B:281:MET:HE2	1.85	0.58
1:A:274:LEU:O	1:A:278:LEU:HB2	2.04	0.57
3:T:86:PHE:CD1	3:T:87:GLU:N	2.73	0.57
1:G:266:LYS:HD3	1:H:267:TYR:CZ	2.40	0.57
3:T:71:GLN:C	3:T:73:LYS:H	2.07	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:LEU:HD11	2:I:4:ILE:HD12	1.86	0.56
3:T:71:GLN:C	3:T:73:LYS:N	2.59	0.56
1:C:270:ILE:HD13	1:C:270:ILE:O	2.06	0.55
1:E:245:GLU:HA	1:E:245:GLU:OE2	2.04	0.55
1:C:266:LYS:HD2	1:D:267:TYR:CE1	2.41	0.55
2:I:0:SER:HB3	2:I:3:ALA:HB3	1.86	0.55
2:I:12:LYS:O	2:I:16:GLU:HG2	2.07	0.55
1:A:273:GLU:OE1	2:I:4:ILE:HG21	2.07	0.55
3:T:86:PHE:O	3:T:90:ARG:HB3	2.06	0.54
1:G:239:ASN:ND2	1:H:239:ASN:HD22	2.05	0.54
1:B:281:MET:HA	1:B:281:MET:HE3	1.89	0.54
3:T:92:GLU:O	3:T:96:LEU:CG	2.50	0.52
3:T:84:SER:CA	3:T:86:PHE:HE2	2.06	0.52
1:B:233:GLY:HA2	1:C:279:ASN:O	2.10	0.52
2:I:12:LYS:NZ	3:T:80:ALA:O	2.41	0.51
1:A:279:ASN:C	1:A:281:MET:N	2.63	0.50
2:I:15:LYS:O	2:I:18:ALA:HB3	2.12	0.50
1:G:267:TYR:O	1:G:270:ILE:HG22	2.11	0.50
1:G:280:ASP:O	1:G:282:THR:HG22	2.11	0.50
2:I:20:ASP:C	2:I:23:GLU:HG2	2.31	0.49
2:I:11:LEU:O	2:I:15:LYS:HG2	2.11	0.49
1:A:279:ASN:O	1:A:281:MET:N	2.39	0.49
3:T:83:ASP:O	3:T:86:PHE:CE2	2.65	0.49
3:T:95:GLU:O	3:T:99:LEU:N	2.45	0.48
3:T:90:ARG:O	3:T:94:GLU:HG3	2.14	0.48
3:T:73:LYS:O	3:T:77:GLU:N	2.43	0.48
3:T:83:ASP:C	3:T:86:PHE:CE2	2.86	0.48
1:B:278:LEU:HD11	2:I:4:ILE:HG23	1.96	0.48
1:B:281:MET:SD	2:I:11:LEU:HG	2.54	0.48
3:T:69:LYS:O	3:T:72:ASN:HB2	2.14	0.47
3:T:86:PHE:CG	3:T:87:GLU:N	2.83	0.46
3:T:73:LYS:HE3	3:T:77:GLU:OE2	2.16	0.46
3:T:82:ILE:C	3:T:84:SER:H	2.19	0.46
1:H:234:GLU:CD	1:H:234:GLU:N	2.66	0.46
3:T:73:LYS:O	3:T:77:GLU:CB	2.61	0.46
1:A:276:HIS:HB2	3:T:86:PHE:HB2	1.98	0.46
1:G:239:ASN:HD21	1:H:239:ASN:HD22	1.64	0.46
1:G:260:LEU:HD23	1:G:260:LEU:HA	1.84	0.46
1:A:276:HIS:HB2	3:T:86:PHE:CB	2.46	0.45
1:A:236:LEU:HD13	1:A:236:LEU:HA	1.55	0.45
2:I:31:ARG:O	2:I:34:LYS:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:LEU:HD12	1:A:278:LEU:HA	1.75	0.45
1:A:245:GLU:O	1:A:246:VAL:C	2.55	0.44
1:E:274:LEU:HA	1:F:274:LEU:HD23	1.99	0.44
3:T:98:ALA:O	3:T:99:LEU:C	2.56	0.44
3:T:67:GLN:H	3:T:69:LYS:H	1.65	0.44
1:E:274:LEU:O	1:E:278:LEU:HB2	2.18	0.44
1:G:257:GLU:CG	1:H:256:LEU:HD21	2.48	0.44
1:A:235:LEU:HD23	1:B:235:LEU:HD13	2.00	0.44
3:T:69:LYS:O	3:T:72:ASN:CB	2.67	0.43
2:I:24:GLN:HG3	2:I:27:ASN:HD22	1.84	0.43
2:I:2:ASP:C	2:I:4:ILE:N	2.71	0.43
3:T:67:GLN:HE21	3:T:67:GLN:HB2	1.58	0.43
1:D:252:LEU:HD12	1:D:252:LEU:HA	1.75	0.43
1:A:272:GLU:HG2	3:T:89:ARG:CZ	2.49	0.42
2:I:28:GLU:O	2:I:32:LEU:CB	2.51	0.42
3:T:94:GLU:O	3:T:97:VAL:HB	2.20	0.41
1:F:281:MET:C	1:F:283:SER:N	2.73	0.41
3:T:82:ILE:O	3:T:85:HIS:HB2	2.20	0.41
1:F:233:GLY:HA2	1:G:279:ASN:O	2.21	0.41
2:I:17:ASN:HD22	2:I:17:ASN:HA	1.65	0.41
2:I:27:ASN:O	2:I:31:ARG:HG2	2.20	0.41
1:A:272:GLU:HG2	3:T:89:ARG:NH2	2.36	0.41
1:G:234:GLU:HB3	1:G:238:LYS:NZ	2.36	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:9:GLN:NE2	3:T:84:SER:O[7_455]	1.87	0.33
2:I:9:GLN:CD	3:T:84:SER:O[7_455]	2.11	0.09

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	47/52 (90%)	46 (98%)	0	1 (2%)	7	26
1	B	49/52 (94%)	46 (94%)	3 (6%)	0	100	100
1	C	50/52 (96%)	48 (96%)	2 (4%)	0	100	100
1	D	47/52 (90%)	47 (100%)	0	0	100	100
1	E	47/52 (90%)	46 (98%)	1 (2%)	0	100	100
1	F	49/52 (94%)	46 (94%)	3 (6%)	0	100	100
1	G	50/52 (96%)	48 (96%)	2 (4%)	0	100	100
1	H	47/52 (90%)	47 (100%)	0	0	100	100
2	I	37/40 (92%)	32 (86%)	4 (11%)	1 (3%)	5	19
3	T	32/55 (58%)	26 (81%)	3 (9%)	3 (9%)	0	1
All	All	455/511 (89%)	432 (95%)	18 (4%)	5 (1%)	14	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	280	ASP
2	I	-1	ALA
3	T	70	ARG
3	T	67	GLN
3	T	72	ASN

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	44/47 (94%)	41 (93%)	3 (7%)	16	42
1	B	46/47 (98%)	40 (87%)	6 (13%)	4	12
1	C	47/47 (100%)	38 (81%)	9 (19%)	1	4
1	D	44/47 (94%)	40 (91%)	4 (9%)	9	28
1	E	44/47 (94%)	40 (91%)	4 (9%)	9	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	46/47 (98%)	41 (89%)	5 (11%)	6	19
1	G	47/47 (100%)	40 (85%)	7 (15%)	3	9
1	H	44/47 (94%)	40 (91%)	4 (9%)	9	28
2	I	32/33 (97%)	24 (75%)	8 (25%)	0	2
3	T	31/49 (63%)	17 (55%)	14 (45%)	0	0
All	All	425/458 (93%)	361 (85%)	64 (15%)	3	9

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	LEU
1	A	260	LEU
1	A	268	LYS
1	B	235	LEU
1	B	249	LEU
1	B	252	LEU
1	B	272	GLU
1	B	281	MET
1	B	283	SER
1	C	234	GLU
1	C	249	LEU
1	C	259	GLU
1	C	260	LEU
1	C	270	ILE
1	C	272	GLU
1	C	274	LEU
1	C	282	THR
1	C	283	SER
1	D	236	LEU
1	D	243	GLU
1	D	270	ILE
1	D	274	LEU
1	E	236	LEU
1	E	245	GLU
1	E	260	LEU
1	E	274	LEU
1	F	235	LEU
1	F	249	LEU
1	F	251	LYS
1	F	252	LEU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	F	272	GLU
1	G	234	GLU
1	G	249	LEU
1	G	259	GLU
1	G	260	LEU
1	G	274	LEU
1	G	281	MET
1	G	284	ILE
1	H	234	GLU
1	H	236	LEU
1	H	270	ILE
1	H	274	LEU
2	I	6	LYS
2	I	11	LEU
2	I	15	LYS
2	I	17	ASN
2	I	20	ASP
2	I	31	ARG
2	I	32	LEU
2	I	35	LEU
3	T	66	ILE
3	T	67	GLN
3	T	68	LYS
3	T	69	LYS
3	T	70	ARG
3	T	72	ASN
3	T	75	LEU
3	T	79	GLN
3	T	81	LEU
3	T	84	SER
3	T	87	GLU
3	T	91	LYS
3	T	92	GLU
3	T	99	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	241	HIS
1	A	276	HIS
1	C	276	HIS
1	G	239	ASN

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Mol	Chain	Res	Type
1	H	244	ASN
2	I	17	ASN
2	I	27	ASN
3	T	67	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å <sup>2</sup> )	Q < 0.9
1	A	49/52 (94%)	-0.25	1 (2%) 65 63	20, 37, 88, 102	0
1	B	51/52 (98%)	-0.20	1 (1%) 65 63	22, 32, 58, 119	0
1	C	52/52 (100%)	-0.08	1 (1%) 66 65	34, 46, 77, 98	0
1	D	49/52 (94%)	0.23	3 (6%) 21 17	37, 49, 80, 111	0
1	E	49/52 (94%)	-0.12	0 100 100	24, 38, 69, 95	0
1	F	51/52 (98%)	-0.03	2 (3%) 39 35	26, 35, 58, 109	0
1	G	52/52 (100%)	0.11	2 (3%) 40 36	32, 47, 75, 110	0
1	H	49/52 (94%)	-0.05	1 (2%) 65 63	37, 48, 76, 100	0
2	I	39/40 (97%)	4.26	36 (92%) 0 0	133, 167, 187, 189	2 (5%)
3	T	34/55 (61%)	3.33	24 (70%) 0 0	167, 199, 214, 216	10 (29%)
All	All	475/511 (92%)	0.55	71 (14%) 2 1	20, 45, 186, 216	12 (2%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	25	LEU	11.9
2	I	35	LEU	10.6
2	I	32	LEU	7.5
2	I	22	ALA	6.5
3	T	72	ASN	6.3
2	I	29	VAL	6.0
2	I	7	LYS	5.8
2	I	14	ASP	5.6
1	D	281	MET	5.6
2	I	30	ALA	5.6
2	I	23	GLU	5.5
3	T	86	PHE	5.5
2	I	21	ARG	5.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
3	T	88	ALA	5.4
3	T	75	LEU	5.4
3	T	77	GLU	5.3
3	T	94	GLU	5.3
2	I	8	MET	5.1
2	I	17	ASN	5.0
3	T	99	LEU	4.9
2	I	1	MET	4.9
2	I	6	LYS	4.8
3	T	70	ARG	4.7
3	T	79	GLN	4.7
2	I	28	GLU	4.7
2	I	9	GLN	4.6
2	I	0	SER	4.4
3	T	78	LEU	4.4
3	T	68	LYS	4.3
1	D	280	ASP	4.2
3	T	84	SER	4.1
3	T	76	ILE	4.1
2	I	3	ALA	4.0
2	I	34	LYS	3.9
2	I	18	ALA	3.8
2	I	20	ASP	3.7
3	T	73	LYS	3.7
2	I	27	ASN	3.7
3	T	93	GLU	3.7
2	I	-3	GLY	3.6
3	T	74	ASP	3.6
1	F	282	THR	3.6
3	T	83	ASP	3.6
2	I	11	LEU	3.5
2	I	4	ILE	3.4
1	F	281	MET	3.4
1	G	283	SER	3.2
3	T	87	GLU	3.1
3	T	95	GLU	3.1
2	I	2	ASP	3.0
3	T	96	LEU	2.9
3	T	82	ILE	2.9
2	I	33	LYS	2.9
2	I	26	GLU	2.8
3	T	80	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
2	I	-1	ALA	2.8
2	I	19	LEU	2.7
1	G	282	THR	2.7
2	I	13	LEU	2.5
2	I	5	LYS	2.5
1	B	282	THR	2.5
3	T	69	LYS	2.5
2	I	10	MET	2.5
2	I	-2	ALA	2.4
3	T	71	GLN	2.3
1	H	233	GLY	2.3
1	C	284	ILE	2.2
2	I	31	ARG	2.2
1	D	233	GLY	2.1
2	I	15	LYS	2.1
1	A	276	HIS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.